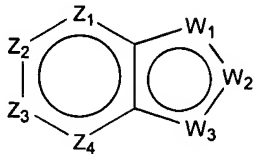
AMENDMENTS TO THE CLAIMS

1. (CURRENTLY AMENDED) A compound of formula (I):



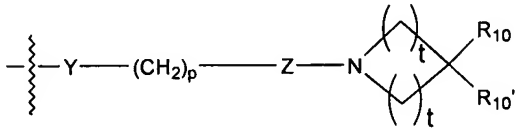
(I)

wherein:

Z_1 is CR_1, Z_2 is CR_2, Z_3 is CR_3, and Z_4 is CR_4;

W_1 is [[O,]] S[[, or NR_5]], W_2 is N or CR_6, and W_3 is CG; ~~W_1 is NG, W_2 is CR_5 or N, and W_3 is CR_6 or N;~~

G is of formula (II):



(II)

Y is O, S, CHOH, -NHC(O)-, -C(O)NH-, -C(O)-, -OC(O)-, -(O)CO-, $-NR_7$-, -CH=N-, or absent;

p is 1, 2, 3, 4 or 5;

Z is CR_8R_9 or absent;

each t is 1, 2, or 3;

each R_1, R_2, R_3, and R_4, independently, is H, amino, hydroxyl, halo, or straight- or branched-chain C_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl, C_{1-6} heteroalkyl, C_{1-6} haloalkyl, -CN, $-CF_3$, $-OR_{11}$, $-COR_{11}$, $-NO_2$, $-SR_{11}$, $-NHC(O)R_{11}$, $-C(O)NR_{12}R_{13}$, $-NR_{12}R_{13}$, $-NR_{11}C(O)NR_{12}R_{13}$, $-SO_2NR_{12}R_{13}$, $-OC(O)R_{11}$, $-O(CH_2)_qNR_{12}R_{13}$, or $-(CH_2)_qNR_{12}R_{13}$, where q is an integer from 2 to 6, or R_1 and R_2 together form -NH-N=N- or R_3 and R_4 together form -NH-N=N-;

each R_5, R_6, and R_7, independently, is H, C_{1-6} alkyl; formyl; C_{3-6} cycloalkyl; C_{5-6} aryl, optionally substituted with halo or C_{1-6} alkyl; or C_{5-6} heteroaryl, optionally substituted with halo or C_{1-6} alkyl;

each R_8 and R_9, independently, is H or straight- or branched-chain C_{1-8} alkyl;

R_{10} is [[H,]] straight- or branched-chain C_{1-8} alkyl, C_{2-8} alkenyl, C_{2-8} alkynyl, C_{1-8} alkylidene, C_{1-8} alkoxy, or C_{1-8} heteroalkyl, ~~C_{1-8} aminoalkyl, C_{1-8} haloalkyl, C_{1-8} alkoxycarbonyl, C_{1-8} hydroxyalkoxy, C_{1-8} hydroxyalkyl, SH, C_{1-8} alkylthio, $-O-CH_2-C_{5-6}$ aryl, $C(O)-C_{5-6}$ aryl~~

~~substituted with C_{1-3} alkyl or halo, C_{5-6} aryl, C_{5-6} cycloalkyl, C_{5-6} heteroaryl, C_{5-6} heterocycloalkyl, $NR_{12}R_{13}$, $-C(O)NR_{12}R_{13}$, $-NR_{11}C(O)NR_{12}R_{13}$, $-CR_{11}R_{12}R_{13}$, $-OC(O)R_{11}$, $-(O)(CH_2)_sNR_{12}R_{13}$ or $-(CH_2)_sNR_{12}R_{13}$, s being an integer from 2 to 8~~;

R_{10}' is H, straight- or branched-chain C_{1-8} alkyl, C_{2-8} alkenyl, C_{2-8} alkynyl, C_{1-8} alkylidene, C_{1-8} alkoxy, C_{1-8} heteroalkyl, C_{1-8} aminoalkyl, C_{1-8} haloalkyl, C_{1-8} alkoxycarbonyl, C_{1-8} hydroxyalkoxy, C_{1-8} hydroxyalkyl, or C_{1-8} alkylthio;

each R_{11}, independently, is H, straight- or branched-chain C_{1-8} alkyl, C_{2-8} alkenyl, C_{2-8} alkynyl, C_{2-8} heteroalkyl, C_{2-8} aminoalkyl, C_{2-8} haloalkyl, C_{1-8} alkoxycarbonyl, C_{2-8} hydroxyalkyl, -C(O)-C_{5-6} aryl substituted with C_{1-3} alkyl or halo, C_{5-6} aryl, C_{5-6} heteroaryl, C_{5-6} cycloalkyl, C_{5-6} heterocycloalkyl, $-C(O)NR_{12}R_{13}$, $-CR_5R_{12}R_{13}$, $-(CH_2)_tNR_{12}R_{13}$, t is an integer from 2 to 8; and

each R_{12} and R_{13}, independently, is H, C_{1-6} alkyl; C_{3-6} cycloalkyl; C_{5-6} aryl, optionally substituted with halo or C_{1-6} alkyl; or C_{5-6} heteroaryl, optionally substituted with halo or C_{1-6} alkyl; ~~or R_{12} and R_{13} together form a cyclic structure;~~

or a pharmaceutically acceptable salt, ester or prodrug thereof.

2. (CURRENTLY AMENDED) The compound of claim 1, wherein each t is 2~~and R_{10} is straight- or branched-chain C_{2-8} alkyl, C_{2-8} alkenyl, C_{2-8} alkynyl, C_{1-8} alkylidene, C_{1-8} alkoxy, or C_{1-8} heteroalkyl~~.

3. (ORIGINAL) The compound of claim 2, wherein R_{10} is n-butyl.

4. (CANCELED)

5. (PREVIOUSLY PRESENTED) The compound of claim 2, wherein each R_1, R_2, R_3, and R_4, independently, is H, hydroxyl, halo, C_{1-6} heteroalkyl, CF_3, $-NO_2$, or straight- or branched-chain C_{1-6} alkyl, or R_1 and R_2 together form -NH-N=N- or R_3 and R_4 together form -NH-N=N-.

6. (ORIGINAL) The compound of claim 2, wherein Y is absent or O, p is 0, 1, 2 or 3, and R_8 and R_9 are H.

7. (ORIGINAL) The compound of claim 6, wherein Z is absent, Y is absent and p is 3.

8. (ORIGINAL) The compound of claim 7, wherein R_{10} is n-butyl.

9-16. (CANCELED)

17. (CURRENTLY AMENDED) The compound of claim 1, wherein the compound is:

~~1-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-indole;~~

~~1-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-benzoimidazole;~~

~~3-methyl-1-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-indole;~~

~~5-bromo-1-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-indole;~~

~~3-formyl-1-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-indole;~~

~~7-bromo-1-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-indole;~~

~~3-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-benzo[*d*]isoxazole;~~

~~3-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-indole;~~

~~3-(2-(4-*n*-butylpiperidine)-1-yl-ethyl)-1*H*-indole;~~

~~3-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-indazole;~~

~~3-(2-(4-n-butylpiperidine)-ethoxy)-7-methyl-benzo[*d*]isoxazole;~~

~~1-(3-(4-methylpiperidine)-1-yl-propyl)-1H-indazole;~~

~~1-(3-(4-pentylpiperidine)-1-yl-propyl)-1H-indazole;~~

~~1-(3-(4-propylpiperidine)-1-yl-propyl)-1H-indazole;~~

~~1-(3-(4-(3-methyl-butyl)-piperidine)-1-yl-propyl)-1H-indazole~~

~~1-(3-(4-pentylidene-piperidine)-1-yl-propyl)-1H-indazole;~~

~~1-(3-(4-propylidene-piperidine)-1-yl-propyl)-1H-indazole~~

1-benzo[*b*]thiophen-2-yl-4-(4-butylpiperidin-1-yl)-butan-1-one

~~4-(4-butylpiperidin-1-yl)-1-(3-methyl-benzofuran-2-yl)-butan-1-one;~~

4-(4-butylpiperidin-1-yl)-1-(5-fluoro-3-methyl-benzo[*b*]thiophen-2-yl)-butan-1-one;

~~1-benzofuran-2-yl-4-(4-butylpiperidin-1-yl)-butan-1-one;~~

1-(3-bromo-benzo[*b*]thiophen-2-yl)-4-(4-butylpiperidin-1-yl)-butan-1-one

1-(3-benzo[*b*]thiophen-2-yl-propyl)-4-butylpiperidine;

~~1-(3-benzofuran-2-yl-propyl)-4-butylpiperidine;~~

~~4-butyl-1-[3-(3-methyl-benzofuran-2-yl)-propyl]-piperidine;~~

4-butyl-1-[3-(5-fluoro-3-methyl-benzo[*b*]thiophen-2-yl)-propyl]-piperidine;

~~2-(3-iodo-propyl)-benzo[*b*]thiophene;~~

1-(3-benzo[*b*]thiophen-2-yl-propyl)-4-methylpiperidine

1-(3-benzo[*b*]thiophen-2-yl-propyl)-4-benzylpiperidine; or

~~1-(3-benzo[*b*]thiophen-2-yl-propyl)-4-(2-methoxy-phenyl)-piperidine;~~

~~1-[3-(4-butylpiperidin-1-yl)-propyl]-1H-benzotriazole;~~

~~1-[3-(4-butylpiperidin-1-yl)-propyl]-1*H*-indole-3-carbaldehyde;~~

~~{1-[3-(4-butylpiperidin-1-yl)-propyl]-1H-indol-3-yl}-methanol;~~

~~1-[3-(4-butylpiperidin-1-yl)-propyl]-2-phenyl-1*H*-benzoimidazole;~~

~~1-[3-(4-butylpiperidin-1-yl)-propyl]-3-chloro-1*H*-indazole;~~

~~1-[3-(4-butylpiperidin-1-yl)-propyl]-6-nitro-1*H*-indazole;~~

~~3-[2-(4-butylpiperidin-1-yl)-ethoxy]-benzo[*d*]isoxazol;~~

~~3-[3-(4-butyl-piperidin-1-yl)-propyl]-1*H*-indole hydrochloride;~~

~~1H-indazole-3-carboxylic acid (2-(4-butylpiperidin)-1-yl-ethyl)-amide;~~

~~1-[3-(4-butylpiperidin-1-yl)-propyl]-5-nitro-1H-indazole;~~

~~1-[3-(4-butyl-piperidin-1-yl)-propyl]-2-methyl-1*H*-indole;~~

~~1-{1-[3-(4-butyl-piperidin-1-yl)-propyl]-1H-indol-3-yl}-ethanone;~~

~~{1-[3-(4-butyl-piperidin-1-yl)-propyl]-1H-indol-3-yl}-acetonitrile;~~

~~1-[3-(4-butyl-piperidin-1-yl)-propyl]-1*H*-indole-3-carbonitrile;~~

~~1-[3-(4-butyl-piperidin-1-yl)-propyl]-5,6-dimethyl-1*H*-benzoimidazole;~~

~~1-[3-(4-butyl-piperidin-1-yl)-propyl]-5(6)-dimethyl-1*H*-benzoimidazole;~~

~~1-[3-(4-butyl-piperidin-1-yl)-propyl]-5-methoxy-1*H*-benzoimidazole;~~

~~{1-[3-(4-butyl-piperidin-1-yl)-propyl]-1H-benzoimidazol-2-yl}-methanol;~~

~~1-[3-(4-butyl-piperidin-1-yl)-propyl]-2-trifuoromethyl-1*H*-benzoimidazole;~~

~~3-[3-(4-butyl-piperidine-1-yl)-propyl]-1H-indazole, HCl;~~

~~3-[3-(4-butyl-piperidine-1-yl)-propyl]-5-nitro-1H-indazole;~~

~~3-[3-(4-butyl-piperidine-1-yl)-propyl]-5,7-dinitro-1H-indazole;~~

3-[3-(4-butyl-piperidin-1-yl)-propyl]-benzo[*d*]isothiazole[[;]].

~~3-[3-(4-butyl-piperidin-1-yl)-propyl]-5-methoxy-1H-indazole;~~

~~3-[3-(4-butyl-piperidin-1-yl)-propyl]-4-methoxy-1H-indazole~~

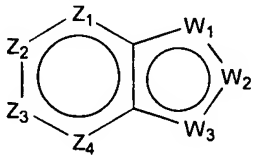
~~3-[3-(4-butyl-piperidin-1-yl)-propyl]-6-methoxy-1H-indazole;~~

~~3-[3-(4-butyl-piperidin-1-yl)-propyl]-1H-indazole-4-ol;~~

~~3-[3-(4-butyl-piperidin-1-yl)-propyl]-1H-indazole-6-ol; or~~

~~3-[3-(4-butyl-piperidin-1-yl)-propyl]-1H-indazole-5-ol.~~

18. (CURRENTLY AMENDED) A pharmaceutical composition comprising an effective amount of a compound of formula (I):



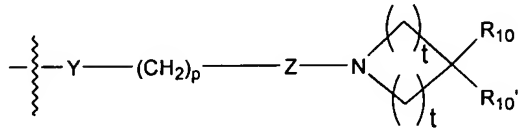
(I)

wherein:

Z_1 is CR_1, Z_2 is CR_2, Z_3 is CR_3, and Z_4 is CR_4;

W_1 is [[O,]] S[[, or NR_5]], W_2 is N or CR_6, and W_3 is CG; ~~W_1 is NG, W_2 is CR_5 or N, and W_3 is CR_6 or N;~~

G is of formula (II):



(II)

Y is O, S, CHOH, -NHC(O)-, -C(O)NH-, -C(O)-, -OC(O)-, -(O)CO-, $-NR_7$-, -CH=N-, or absent;

p is 1, 2, 3, 4 or 5;

Z is CR_8R_9 or absent;

each t is 1, 2, or 3;

each R_1, R_2, R_3, and R_4, independently, is H, amino, hydroxyl, halo, or straight- or branched-chain C_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl, C_{1-6} heteroalkyl, C_{1-6} haloalkyl, -CN, $-CF_3$, $-OR_{11}$, $-COR_{11}$, $-NO_2$, $-SR_{11}$, $-NHC(O)R_{11}$, $-C(O)NR_{12}R_{13}$, $-NR_{12}R_{13}$, $-NR_{11}C(O)NR_{12}R_{13}$, $-SO_2NR_{12}R_{13}$, $-OC(O)R_{11}$, $-O(CH_2)_qNR_{12}R_{13}$, or $-(CH_2)_qNR_{12}R_{13}$, where q is an integer from 2 to 6, or R_1 and R_2 together form -NH-N=N- or R_3 and R_4 together form -NH-N=N-;

each R_5, R_6, and R_7, independently, is H, C_{1-6} alkyl; formyl; C_{3-6} cycloalkyl; C_{5-6} aryl, optionally substituted with halo or C_{1-6} alkyl; or C_{5-6} heteroaryl, optionally substituted with halo or C_{1-6} alkyl;

each R_8 and R_9, independently, is H or straight- or branched-chain C_{1-8} alkyl;

R_{10} is [[H,]] straight- or branched-chain C_{1-8} alkyl, C_{2-8} alkenyl, C_{2-8} alkynyl, C_{1-8} alkylidene, C_{1-8} alkoxy, or C_{1-8} heteroalkyl~~, C_{1-8} aminoalkyl, C_{1-8} haloalkyl, C_{1-8} alkoxycarbonyl,~~

~~C_{1-8} hydroxyalkoxy, C_{1-8} hydroxyalkyl, SH, C_{1-8} alkylthio, $-O-CH_2-C_{5-6}$ aryl, $-C(O)-C_{5-6}$ aryl substituted with C_{1-3} alkyl or halo, C_{5-6} aryl, C_{5-6} cycloalkyl, C_{5-6} heteroaryl, C_{5-6} heterocycloalkyl, $NR_{12}R_{13}$, $C(O)NR_{12}R_{13}$, $-NR_{11}C(O)NR_{12}R_{13}$, $-CR_{11}R_{12}R_{13}$, $-OC(O)R_{11}$, $-(O)(CH_2)_sNR_{12}R_{13}$ or $-(CH_2)_sNR_{12}R_{13}$, s being an integer from 2 to 8~~;

R_{10}' is H, straight- or branched-chain C_{1-8} alkyl, C_{2-8} alkenyl, C_{2-8} alkynyl, C_{1-8} alkylidene, C_{1-8} alkoxy, C_{1-8} heteroalkyl, C_{1-8} aminoalkyl, C_{1-8} haloalkyl, C_{1-8} alkoxycarbonyl, C_{1-8} hydroxyalkoxy, C_{1-8} hydroxyalkyl, or C_{1-8} alkylthio;

each R_{11}, independently, is H, straight- or branched-chain C_{1-8} alkyl, C_{2-8} alkenyl, C_{2-8} alkynyl, C_{2-8} heteroalkyl, C_{2-8} aminoalkyl, C_{2-8} haloalkyl, C_{1-8} alkoxycarbonyl, C_{2-8} hydroxyalkyl, -C(O)-C_{5-6} aryl substituted with C_{1-3} alkyl or halo, C_{5-6} aryl, C_{5-6} heteroaryl, C_{5-6} cycloalkyl, C_{5-6} heterocycloalkyl, $-C(O)NR_{12}R_{13}$, $-CR_5R_{12}R_{13}$, $-(CH_2)_tNR_{12}R_{13}$, t is an integer from 2 to 8; and

each R_{12} and R_{13}, independently, is H, C_{1-6} alkyl; C_{3-6} cycloalkyl; C_{5-6} aryl, optionally substituted with halo or C_{1-6} alkyl; or C_{5-6} heteroaryl, optionally substituted with halo or C_{1-6} alkyl; ~~or R_{12} and R_{13} together form a cyclic structure;~~

or a pharmaceutically acceptable salt, ester or prodrug thereof.

19. (CURRENTLY AMENDED) A pharmaceutical composition of Claim 18, wherein each t is 2~~ and R_{10} is straight- or branched-chain C_{2-8} alkyl, C_{2-8} alkenyl, C_{2-8} alkynyl, C_{1-8} alkylidene, C_{1-8} alkoxy, or C_{1-8} heteroalkyl~~.

20. (ORIGINAL) A pharmaceutical composition of Claim 19, wherein R_{10} is n-butyl.

21. (CANCELED)

22. (PREVIOUSLY PRESENTED) A pharmaceutical composition of Claim 19, wherein each R_1, R_2, R_3, and R_4, independently, is H, hydroxyl, halo, C_{1-6}heteroalkyl, CF_3, $-NO_2$, or straight- or branched-chain C_{1-6} alkyl, or R_1 and R_2 together form -NH-N=N- or R_3 and R_4 together form -NH-N=N-.

23. (ORIGINAL) A pharmaceutical composition of Claim 19, wherein Y is absent or O, p is 0, 1, 2 or 3, and R_8 and R_9 are H.

24. (ORIGINAL) A pharmaceutical composition of Claim 23, wherein Z is absent, Y is absent and p is 3.

25. (ORIGINAL) A pharmaceutical composition of Claim 24, wherein R_{10} is n-butyl.

26-33. (CANCELED)

34. (CURRENTLY AMENDED) A pharmaceutical composition of Claim 19, wherein the compound is:

~~1-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-indole;~~

~~1-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-benzoimidazole;~~

~~3-methyl-1-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-indole;~~

~~5-bromo-1-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-indole;~~

~~3-formyl-1-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-indole;~~

~~7-bromo-1-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-indole;~~

~~3-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-benzo[*d*]isoxazole;~~

~~3-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-indole;~~

~~3-(2-(4-*n*-butylpiperidine)-1-yl-ethyl)-1*H*-indole;~~

~~3-(3-(4-*n*-butylpiperidine)-1-yl-propyl)-1*H*-indazole;~~

~~3-(2-(4-n-butylpiperidine)-ethoxy)-7-methyl-benzo[*d*]isoxazole;~~

~~1-(3-(4-methylpiperidine)-1-yl-propyl)-1H-indazole;~~

~~1-(3-(4-pentylpiperidine)-1-yl-propyl)-1H-indazole;~~

~~1-(3-(4-propylpiperidine)-1-yl-propyl)-1H-indazole;~~

~~1-(3-(4-(3-methyl-butyl)-piperidine)-1-yl-propyl)-1H-indazole~~

~~1-(3-(4-pentylidene-piperidine)-1-yl-propyl)-1H-indazole;~~

~~1-(3-(4-propylidene-piperidine)-1-yl-propyl)-1H-indazole~~

1-benzo[*b*]thiophen-2-yl-4-(4-butylpiperidin-1-yl)-butan-1-one

~~4-(4-butylpiperidin-1-yl)-1-(3-methyl-benzofuran-2-yl)-butan-1-one;~~

4-(4-butylpiperidin-1-yl)-1-(5-fluoro-3-methyl-benzo[*b*]thiophen-2-yl)-butan-1-one;

~~1-benzofuran-2-yl-4-(4-butylpiperidin-1-yl)-butan-1-one;~~

1-(3-bromo-benzo[*b*]thiophen-2-yl)-4-(4-butylpiperidin-1-yl)-butan-1-one

1-(3-benzo[*b*]thiophen-2-yl-propyl)-4-butylpiperidine;

~~1-(3-benzofuran-2-yl-propyl)-4-butylpiperidine;~~

~~4-butyl-1-[3-(3-methyl-benzofuran-2-yl)-propyl]-piperidine;~~

4-butyl-1-[3-(5-fluoro-3-methyl-benzo[*b*]thiophen-2-yl)-propyl]-piperidine;

~~2-(3-iodo-propyl)-benzo[*b*]thiophene;~~

1-(3-benzo[*b*]thiophen-2-yl-propyl)-4-methylpiperidine

1-(3-benzo[*b*]thiophen-2-yl-propyl)-4-benzylpiperidine; or

~~1-(3-benzo[*b*]thiophen-2-yl-propyl)-4-(2-methoxy-phenyl)-piperidine;~~

~~1-[3-(4-butylpiperidin-1-yl)-propyl]-1H-benzotriazole;~~

~~1-[3-(4-butylpiperidin-1-yl)-propyl]-1*H*-indole-3-carbaldehyde;~~

~~{1-[3-(4-butylpiperidin-1-yl)-propyl]-1H-indol-3-yl}-methanol;~~

~~1-[3-(4-butylpiperidin-1-yl)-propyl]-2-phenyl-1*H*-benzoimidazole;~~

~~1-[3-(4-butylpiperidin-1-yl)-propyl]-3-chloro-1*H*-indazole;~~

~~1-[3-(4-butylpiperidin-1-yl)-propyl]-6-nitro-1*H*-indazole;~~

~~3-[2-(4-butylpiperidin-1-yl)-ethoxy]-benzo[*d*]isoxazol;~~

~~3-[3-(4-butyl-piperidin-1-yl)-propyl]-1*H*-indole hydrochloride;~~

~~1H-indazole-3-carboxylic acid (2-(4-butylpiperidin)-1-yl-ethyl)-amide;~~

~~1-[3-(4-butylpiperidin-1-yl)-propyl]-5-nitro-1H-indazole;~~

~~1-[3-(4-butyl-piperidin-1-yl)-propyl]-2-methyl-1*H*-indole;~~

~~1-{1-[3-(4-butyl-piperidin-1-yl)-propyl]-1H-indol-3-yl}-ethanone;~~

~~{1-[3-(4-butyl-piperidin-1-yl)-propyl]-1H-indol-3-yl}-acetonitrile;~~

~~1-[3-(4-butyl-piperidin-1-yl)-propyl]-1*H*-indole-3-carbonitrile;~~

~~1-[3-(4-butyl-piperidin-1-yl)-propyl]-5,6-dimethyl-1*H*-benzoimidazole;~~

~~1-[3-(4-butyl-piperidin-1-yl)-propyl]-5(6)-dimethyl-1*H*-benzoimidazole;~~

~~1-[3-(4-butyl-piperidin-1-yl)-propyl]-5-methoxy-1*H*-benzoimidazole;~~

~~{1-[3-(4-butyl-piperidin-1-yl)-propyl]-1H-benzoimidazol-2-yl}-methanol;~~

~~1-[3-(4-butyl-piperidin-1-yl)-propyl]-2-trifuoromethyl-1*H*-benzoimidazole;~~

~~3-[3-(4-butyl-piperidine-1-yl)-propyl]-1H-indazole, HCl;~~

~~3-[3-(4-butyl-piperidine-1-yl)-propyl]-5-nitro-1H-indazole;~~

~~3-[3-(4-butyl-piperidine-1-yl)-propyl]-5,7-dinitro-1H-indazole;~~

3-[3-(4-butyl-piperidin-1-yl)-propyl]-benzo[*d*]isothiazole[[;]].

~~3-[3-(4-butyl-piperidin-1-yl)-propyl]-5-methoxy-1H-indazole;~~

~~3-[3-(4-butyl-piperidin-1-yl)-propyl]-4-methoxy-1H-indazole~~

~~3-[3-(4-butyl-piperidin-1-yl)-propyl]-6-methoxy-1H-indazole;~~

~~3-[3-(4-butyl-piperidin-1-yl)-propyl]-1H-indazole-4-ol;~~

~~3-[3-(4-butyl-piperidin-1-yl)-propyl]-1H-indazole-6-ol; or~~

~~3-[3-(4-butyl-piperidin-1-yl)-propyl]-1H-indazole-5-ol.~~

35-76. (CANCELED)